92197 SEARCH REQUEST FORM

Access DB#

Scientific and Technical Information Center

Mail Box and Bldg/Room Loo	CYER	Results Format Preferred (circle): PAPER DISK E-MAIL										
If more than one search is submitted, please prioritize searches in order of need.												
Please provide a detailed statement of the search topic, and describe as specifically as possible the subject matter to be searched. Include the elected species or structures, keywords, synonyms, acronyms, and registry numbers, and combine with the concept or utility of the invention. Define any terms that may have a special meaning. Give examples or relevant citations, authors, etc., if known. Please attach a copy of the cover sheet, pertinent claims, and abstract.												
Title of Invention:												
Inventors (please provide full name	es):											
Earliest Priority Filing Date:												
Appropriate serial number.	Mary Jane Ruhl 1. Info. Specialist, STIC TC-1600 CM-1, Room 6A-06 Phone: 605-1155	A=H/Ab NO PY 2002 a 2003										
STAFF USE ONLY	Type of Search NA Sequence (#)	Vendors and cost where applicable										
Searcher Phone #:	AA Sequence (#)											
Searcher Location:	<u> </u>	Questel/Orbit										
Date Searcher Picked Up:	Bibliographic											
Date Completed: 4/24/0-3	Litigation	Lexis/Nexis										
Searcher Prep & Review Time:	Fulltext	Sequence Systems										
Clerical Prep Time:	Patent Family	WWW/internet										
Online Time:	Other	Other (enecify)										

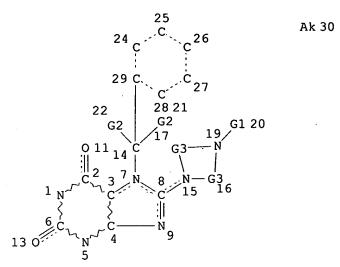
PTO-1590 (8-01)

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(FILE 'HOME' ENTERED AT 10:12:18 ON 24 APR 2003)
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FILE 'REGISTRY' ENTERED AT 10:12:32 ON 24 APR 2003
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L1
L2
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L19
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L20
                   STR L19
      2 S L20 FULL 2 compds in Reg. -pls see done stat for structure
FILE 'HCAPLUS' ENTERED AT 10:56:37 ON 24 APR 2003

1 S L22 / cit in CA Plus - its 2002, attacked
                 1 S L20
L21
L22
L23
                         (Ocits prior to 2002)
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VAR G1=H/CH3
VAR G2=C/H
REP G3=(1-7) CH2
NODE ATTRIBUTES:
DEFAULT MLEVEL IS ATOM
GGCAT IS LOC UNS AT 30
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES: RING(S) ARE ISOLATED OR EMBEDDED NUMBER OF NODES IS 26

STEREO ATTRIBUTES: NONE

L22 2 SEA FILE=REGISTRY SSS FUL L20 L23 1 SEA FILE=HCAPLUS ABB=ON L22

=> d ibib abs hitstr 1-1

L23 ANSWER 1 OF 1 HCAPLUS COPYRIGHT 2003 ACS ACCESSION NUMBER: 2002:31449 HCAPLUS

DOCUMENT NUMBER: 136:102402

TITLE: Preparation of piperazinylpurinediones as inhibitors

of dipeptidylpeptidase IV.

INVENTOR(S): Kanstrup, Anders Bendtz; Christiansen, Lise Brown;

Lundbeck, Jane Marie; Sams, Christian K.; Kristiansen,

Marit

PATENT ASSIGNEE(S): Novo Nordisk A/S, Den.

SOURCE: PCT Int. Appl., 135 pp.

CODEN: PIXXD2
DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

	PATENT NO.				KIND DATE				APPLICATION NO.						DATE				
	WO	vo 2002002560				A2 20020110				- W	0 20	01-D	20010704						
		W:	ΑE,	AG,	AL,	AM,	AT,	AU,	AZ,	BA,	BB,	BG,	BR,	BY,	ΒZ,	CA,	CH,	CN,	
			CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	ES,	FI,	GB,	GD,	GE,	GH,	
			GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	ΚE,	KG,	KΡ,	KR,	ΚZ,	LC,	LK,	LR,	
			LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MZ,	NO,	NZ,	PL,	PT,	
			RO,	RU,	SD,	SE,	SG,	SI,	SK,	SL,	TJ,	TM,	TR,	TT,	TZ,	UA,	UG,	US,	
			UZ,	VN,	YU,	ZA,	ZW,	AM,	AZ,	BY,	KG,	KZ,	MD,	RU,	TJ,	TM			
		RW:	GH,	GM,	KE,	LS,	MW,	MZ,	SD,	SL,	SZ,	TZ,	UG,	ZW,	ΑT,	BE,	CH,	CY,	
			DE,	DK,	ES,	FI,	FR,	GB,	GR,	ΙE,	IT,	LU,	MC,	NL,	PT,	SE,	TR,	BF,	
			ВJ,	CF,	CG,	CI,	CM,	GΑ,	GN,	GW,	ML,	MR,	NE,	SN,	TD,	TG			
	ΑU	2001068958			A5 20020114			AU 2001-68958						20010704					
	ΕP	P 1301187			A2 20030416			EP 2001-947211						20010704					
		R:	AT,	ΒĒ,	CH,	DE,	DK,	ES,	FR,	GB,	GR,	IT,	LI,	LU,	NL,	SE,	MC,	PT,	
			IE,	SI,	LT,	LV,	FI,	RO,	MK,	CY,	AL,	TR							
US 2002161001 A1 20021031										US 2001-935149					2001	0822			
PRIOF	RIT	Y APP	LN.	INFO	. :				1	DK 2	-000	1040		Α	2000	0704			
							US 2000-223240P					Ρ	2000	0704					
									1	WO 2	001-	DK46	7	W	2001	0704			
OWURD COUDCE (C). MADDAM 126.102402																			

OTHER SOURCE(S):

MARPAT 136:102402

GI

AB Title compds. [I; m, n = 1, 2; Q = (CH2)n; Q1 = (CH2)m; R1 = CO, CS, (substituted) alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, aralkyl, heteroaryl, heteroarylalkyl; R2 = H, (substituted) alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, aralkyl, heteroarylalkyl, heteroaryl, cyano, halo, OH, NO2, SH, SR5, SOR5, SO2R5,

CO2H, CO2R4, CON(R5)2, alkoxy, alkenyloxy, alkynyloxy, aryloxy, heteroaryloxy; R3 = H, alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, aralkyl, heteroarylalkyl, heteroaryl, alkoxyalkyl, CO2H, cyano, NO2, halo, OH; R3R3C, R4R4C = spiro system; R4 = (substituted) alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, heteroaryl, cyano, halo, OH, NO2, CF3, alkoxy, alkenyloxy, alkynyloxy, aryloxy, heteroaryloxy, etc.; R5 = H, (substituted) alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, arylalkyl, heteroaryl, heteroarylalkyl; R6 = H, (substituted) alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, heteroaryl; R7 = H, (substituted) alkyl, alkenyl, alkynyl, cycloalkyl, cycloheteroalkyl, aryl, heteroaryl; R9 = H, halo, (substituted) alkyl, aryl; R10 = H, halo; R9R10 = atoms to form a cyclopropyl ring], were prepd. as DPP-IV inhibitors (no data). Thus, 1,4-diazepan-6-ylmethanol, 7-benzyl-8-chloro-1,3-dimethyl-3,7dihydropurine-2,6-dione, and K2CO3 were stirred at room temp. for 24 h, heated to 60.degree. for 3 h, heated to 95.degree. for 5 h, and heated to 120.degree. for 2 h to give 7-benzyl-8-(6-hydroxymethyl-1,4-diazepan-1-yl)-1,3-dimethyl-3,7-dihydropurine-2,6-dione.

IT 389060-65-5P 389062-41-3P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(prepn. of piperazinylpurinediones as inhibitors of dipeptidylpeptidase ${\tt IV})$

RN 389060-65-5 HCAPLUS

CN 1H-Purine-2,6-dione, 8-(hexahydro-1H-1,4-diazepin-1-yl)-3,7-dihydro-3-methyl-7-(phenylmethyl)-1-(3-phenyl-2-propenyl)- (9CI) (CA INDEX NAME)

RN 389062-41-3 HCAPLUS

CN 1H-Purine-2,6-dione, 8-(hexahydro-1H-1,4-diazepin-1-yl)-3,7-dihydro-3-methyl-7-(phenylmethyl)-1-(3-phenyl-2-propenyl)-, trifluoroacetate (9CI) (CA INDEX NAME)

CM 1

CRN 389060-65-5 CMF C27 H30 N6 O2

$$\begin{array}{c|c} Me \\ \hline 0 & N & N & NH \\ \hline Ph-CH=CH-CH_2 & CH_2-Ph \\ \hline \end{array}$$

CM 2

CRN 76-05-1 CMF C2 H F3 O2

=> d 16 1-10

ANSWER 1 OF 10 REGISTRY COPYRIGHT 2003 ACS L6

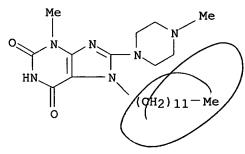
477333-67-8 REGISTRY RN

1H-Purine-2,6-dione, 7-dodecyl-3,7-dihydro-3-methyl-8-(4-methyl-1-CN piperazinyl) - (9CI) (CA INDEX NAME)

FS 3D CONCORD

C23 H40 N6 O2 MF

SR CAS Registry Services



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ANSWER 2 OF 10 REGISTRY COPYRIGHT 2003 ACS L6

454706-71-9 REGISTRY

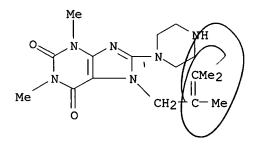
1H-Purine-2,6-dione, 7-(2,3-dimethyl-2-butenyl)-3,7-dihydro-1,3-dimethyl-8-(1-piperazinyl) - (9CI) (CA INDEX NAME)

3D CONCORD

MF C17 H26 N6 O2

SR CA

LC STN Files: CA, CAPLUS



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1957 TO DATE)

1 REFERENCES IN FILE CAPLUS (1957 TO DATE)

L6 ANSWER 3 OF 10 REGISTRY COPYRIGHT 2003 ACS

RN 402470-56-8 REGISTRY

CN 1H-Purine-2,6-dione, 3,7-dihydro-8-[4-(2-hydroxyethyl)-1-piperazinyl]-3methyl-7-nonyl- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C21 H36 N6 O3 SR Chemical Library

LC STN Files: CHEMCATS

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L6 ANSWER 4 OF 10 REGISTRY COPYRIGHT 2003 ACS

RN · 378201-33-3 REGISTRY

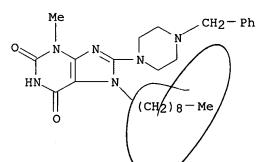
CN 1H-Purine-2,6-dione, 3,7-dihydro-3-methyl-7-nonyl-8-[4-(phenylmethyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C26 H38 N6 O2

SR Chemical Library

LC STN Files: CHEMCATS



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- L6 ANSWER 5 OF 10 REGISTRY COPYRIGHT 2003 ACS
- RN 377060-82-7 REGISTRY
- CN lH-Purine-2,6-dione, 3,7-dihydro-3-methyl-7-octyl-8-[4-(phenylmethyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)
- FS 3D CONCORD
- MF C25 H36 N6 O2
- SR Chemical Library
- LC STN Files: CHEMCATS

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- L6 ANSWER 6 OF 10 REGISTRY COPYRIGHT 2003 ACS
- RN 371927-37-6 REGISTRY
- CN 1H-Purine-2,6-dione, 7-hexyl-3,7-dihydro-3-methyl-8-(1-piperazinyl)- (9CI) (CA INDEX NAME)
- FS 3D CONCORD
- MF C16 H26 N6 O2
- SR Chemical Library
- LC STN Files: CHEMCATS

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- L6 ANSWER 7 OF 10 REGISTRY COPYRIGHT 2003 ACS
- RN 329702-29-6 REGISTRY
- CN 1H-Purine-2,6-dione, 3,7-dihydro-3-methyl-7-nonyl-8-(1-piperazinyl)- (9CI) (CA INDEX NAME)
- FS 3D CONCORD
- MF C19 H32 N6 O2
- SR Chemical Library
- LC STN Files: CHEMCATS

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L6 ANSWER 8 OF 10 REGISTRY COPYRIGHT 2003 ACS

RN 313471-56-6 REGISTRY

CN 1H-Purine-2,6-dione, 7-heptyl-3,7-dihydro-3-methyl-8-(4-methyl-1-piperazinyl)- (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C18 H30 N6 O2

SR Chemical Library

LC STN Files: CHEMCATS

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PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L6 ANSWER 9 OF 10 REGISTRY COPYRIGHT 2003 ACS

RN 304876-71-9 REGISTRY

CN 1H-Purine-2,6-dione, 7-hexadecyl-3,7-dihydro-3-methyl-8-(1-piperazinyl)-(9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C26 H46 N6 O2

SR Chemical Library

LC STN Files: CHEMCATS

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PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L6 ANSWER 10 OF 10 REGISTRY COPYRIGHT 2003 ACS

RN 106939-21-3 REGISTRY

CN 1H-Purine-2,6-dione, 7-heptyl-3,7-dihydro-3-methyl-8-(1-piperazinyl)-(9CI) (CA INDEX NAME)

FS 3D CONCORD

Berch 09/935,149

MF C17 H28 N6 O2

SR CA

LC STN Files: CA, CAPLUS, CASREACT, CHEMCATS, TOXCENTER

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 1 REFERENCES IN FILE CA (1957 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1957 TO DATE)